IN THE CLAIMS

This listing of claims will replace all prior versions, and listings, of claims in the application.

LISTING OF CLAIMS:

1. (Currently Amended) 1. A compound of formula (I):

S^{ag}-T^{ag}

(I)

or a stereoisomer or a pharmaceutically acceptable salt form thereof, wherein:

S^{ag} is ~~R^A or R^B;~~

~~wherein~~

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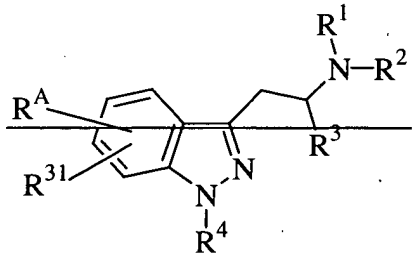
~~R^A is~~ $-NR^{12}R^{13}$;

~~R^B is~~ $-NR^{12}R^{13}$ or aryl substituted with 0-5 R^{33}; and

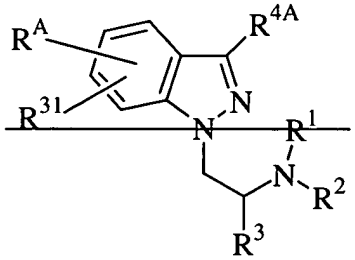
aryl is phenyl, pyridyl, or naphthyl; and

T^{ag} is a heterocyclic serotonin receptor ligand template selected from:

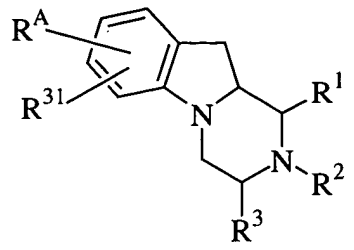
~~1)~~



~~2)~~



3 1)



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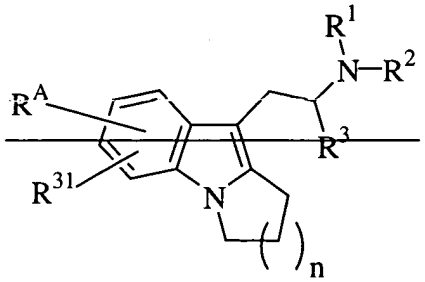

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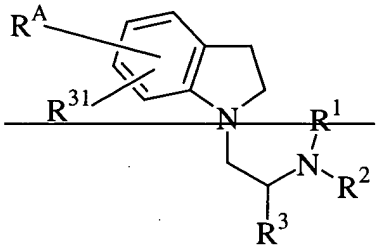

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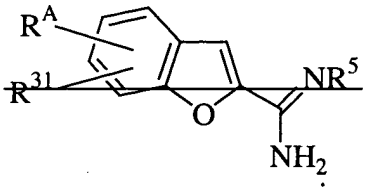

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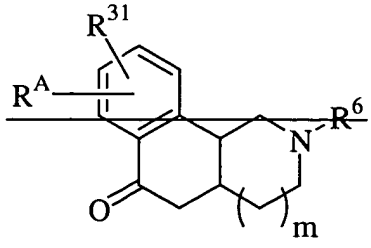

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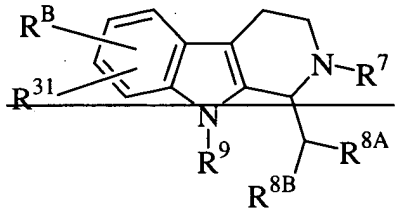

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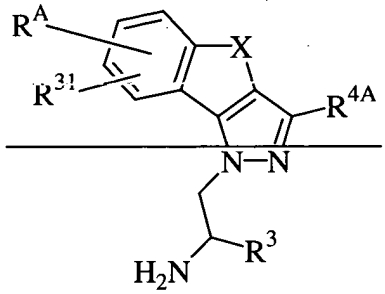

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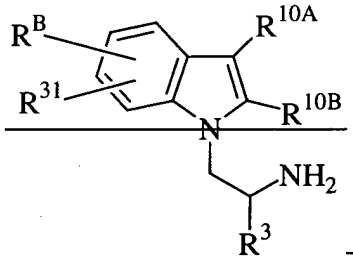
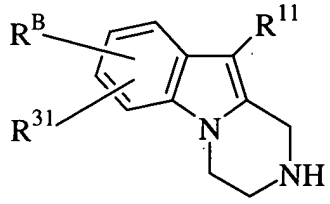
~~10)~~



~~and~~ or

~~11~~ 2)



wherein:

R^A is $-NR^{12}R^{13}$;

R^B is $-NR^{12}R^{13}$ or aryl substituted with 0-5 R^{33}; and aryl is phenyl, pyridyl or naphthyl;

~~X is -CH=CH-, $-CR^1R^2-$, or $-CR^1R^2-CR^1R^2-$;~~

~~n is 1, 2, or 3;~~

~~m is 0 or 1;~~

R^1 is H or C_{1-4} alkyl;

R^2 is H or C_{1-4} alkyl;

R^3 is H or C_{1-4} alkyl;

~~R^4 is H or C_{1-4} alkyl;~~

~~R^{4A} is H, C_{1-4} alkyl or C_{1-4} alkyloxy;~~

~~R^5 is H, -OH or C_{1-4} alkyloxy;~~

~~R^6 is H, C_{1-6} alkyl, C_{2-6} alkenyl, or C_{1-2} alkyl substituted with R^{6A};~~

~~R^{6A} is phenyl, cyclopropyl, cyclobutyl, cyclopentyl or cyclohexyl;~~

~~R^7 is H, C_{1-6} alkyl, C_{2-6} alkenyl, cyclopropyl, cyclobutyl, cyclopentyl or cyclohexyl;~~

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~~R^{8A} is a 5-10 membered heterocyclic ring system containing from 1-3 heteroatoms selected from the group~~

~~benzimidazolyl, benzimidazolinyl,benztriazolyl, benzisoxazolyl, benzisoxazolinyl, benzoxazolyl, benzoxazolinyl, benzthiazolyl, benzisothiazolyl, indolyl, indolinyl, isoindolinyl, indazolyl, isatinoyl, isoxazolopyridinyl, isothiazolopyridinyl, thiazolopyridinyl, oxazolopyridinyl, oxindolyl, oxazolidinyl, imidazolopyridinyl, and pyrazolopyridinyl;~~

~~wherein said heterocyclic ring system is substituted with 0-2 R^{41};~~

~~R^{8B} is H, C_{1-6} alkyl, C_{2-6} alkenyl, cyclopropyl, cyclobutyl, cyclopentyl or cyclohexyl;~~

~~R^9 is H, C_{1-6} alkyl, C_{2-6} alkenyl, cyclopropyl, cyclobutyl, cyclopentyl or cyclohexyl;~~

~~R^{10A} and R^{10b}, at each occurrence, are independently selected from~~

~~H, -OH, halo, $-CF_3$, $-OCF_3$, C_{1-6} alkyl, C_{1-6} alkyloxy-, cyclopropyl, cyclobutyl, cyclopentyl or cyclohexyl;~~

R^{11} is H, C_{1-6} alkoxy or C_{1-6} alkyl-S-;

R^{12} is aryl substituted with 0-5 R^{33};

R^{13} is selected from

H, C_{1-4} alkyl, C_{2-4} alkenyl, and C_{2-4} alkynyl;

alternatively, R^{12} and R^{13} join to form a 5- or 6-membered ring optionally substituted with -O- or -$N(R^{14})$-;

alternatively, R^{12} and R^{13} when attached to N may be combined to form a 9- or 10-membered bicyclic heterocyclic ring system containing from 1-3 heteroatoms selected from the group

consisting of N, O, and S, wherein said bicyclic heterocyclic ring system is unsaturated or partially saturated, wherein said bicyclic heterocyclic ring system is substituted with 0-3 R^{16};

R^{14} is H or C_{1-4} alkyl;

R^{16}, at each occurrence, is independently selected from

H, OH, halo, CN, NO_2, CF_3, SO_2R^{45}, $NR^{46}R^{47}$, -C(=O)H,

C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} haloalkyl,

C_{1-3} haloalkyl-oxy-, and C_{1-3} alkyloxy-;

R^{31}, at each occurrence, is independently selected from

H, -OH, halo, -CN, -NO_2, -CF_3, -OCF_3, C_{1-6} alkyl,

C_{1-4} alkyl-C(=O)-, C_{1-4} alkyloxy-, C_{1-4} alkylthio-,

C_{1-4} alkyl-S(=O)-, and C_{1-4} alkyl-SO_2-;

R^{33}, at each occurrence, is independently selected from

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H, OH, halo, CN, NO_2, CF_3, SO_2R^{45}, $NR^{46}R^{47}$, -C(=O)H, phenyl, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl,

C_{3-6} cycloalkyl, C_{1-4} haloalkyl, C_{1-4} haloalkyl-oxy-, C_{1-4} alkyloxy-, C_{1-4} alkylthio-, C_{1-4} alkyl-C(=O)-,

C_{1-4} alkyl-C(=O)NH-, C_{1-4} alkyl-OC(=O)-,

C_{1-4} alkyl-C(=O)O-, C_{3-6} cycloalkyl-oxy-,

C_{3-6} cycloalkylmethyl-oxy-;

C_{1-6} alkyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, $-SO_2R^{45}$, $-NR^{46}R^{47}$, $NR^{46}R^{47}C(=O)$-, or $(C_{1-4}$ alkyl$)CO_2$-; and

C_{2-6} alkenyl substituted with OH, methoxy, ethoxy, propoxy, butoxy, $-SO_2R^{45}$, $-NR^{46}R^{47}$, $NR^{46}R^{47}C(=O)$-, or $(C_{1-4}$ alkyl$)CO_2$-;

~~R^{41}, at each occurrence, is independently selected from~~

~~H, -OH, F, Cl, -CF_3, -OCF_3, methyl, ethyl, methoxy, and ethoxy;~~

R^{45} is C_{1-4} alkyl;

R^{46}, at each occurrence, is independently selected from H and C_{1-4} alkyl; and

R^{47}, at each occurrence, is independently selected from H, C_{1-4} alkyl, -C(=O)NH(C_{1-4} alkyl), -SO_2(C_{1-4} alkyl),

-C(=O)O(C_{1-4} alkyl), -C(=O)(C_{1-4} alkyl), and -C(=O)H.

2. (Original) A compound of Claim 1 of formula (I), wherein:

R^A is $-NR^{12}R^{13}$;

R^B is $-NR^{12}R^{13}$;

phenyl- substituted with 0-5 fluoro;

naphthyl- substituted with 0-3 R^{33};

2-($H_3CCH_2C(=O)$)-phenyl- substituted with R^{33};

2-($H_3CC(=O)$)-phenyl- substituted with R^{33};

2-(HC(=O))-phenyl- substituted with R^{33};

2-($H_3CCH(OH)$)-phenyl- substituted with R^{33};

2-($H_3CCH_2CH(OH)$)-phenyl- substituted with R^{33};

#4 2-($HOCH_2$)-phenyl- substituted with R^{33};

2-($HOCH_2CH_2$)-phenyl- substituted with R^{33};

2-(H_3COCH_2)-phenyl- substituted with R^{33};

2-($H_3COCH_2CH_2$)-phenyl- substituted with R^{33};

2-($H_3CCH(OMe)$)-phenyl- substituted with R^{33};

2-($H_3COC(=O)$)-phenyl- substituted with R^{33};

2-($HOCH_2CH=CH$)-phenyl- substituted with R^{33};

2-((MeOC=O)CH=CH)-phenyl- substituted with R^{33};

2-(methyl)-phenyl- substituted with R^{33};

2-(ethyl)-phenyl- substituted with R^{33};

2-(i-propyl)-phenyl- substituted with R^{33};

2-(F_3C)-phenyl- substituted with R^{33};

2-(NC)-phenyl- substituted with R^{33};

2-(H_3CO)-phenyl- substituted with R^{33};

2-(fluoro)-phenyl- substituted with R^{33};

2-(chloro)-phenyl- substituted with R^{33};

3-(NC)-phenyl- substituted with R^{33};

3-(H_3CO)-phenyl- substituted with R^{33};

3-(fluoro)-phenyl- substituted with R^{33};

3-(chloro)-phenyl- substituted with R^{33};

4-(NC)-phenyl- substituted with R^{33};

4-(fluoro)-phenyl- substituted with R^{33};

4-(chloro)-phenyl- substituted with R^{33};

4-(H_3CS)-phenyl- substituted with R^{33};

4-(H_3CO)-phenyl- substituted with R^{33};

4-(ethoxy)-phenyl- substituted with R^{33};

4-(i-propoxy)-phenyl- substituted with R^{33};

4-(i-butoxy)-phenyl- substituted with R^{33};

4-($H_3CCH_2CH_2C(=O)$)-phenyl- substituted with R^{33};

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4-($(H_3C)_2CHC(=O)$)-phenyl- substituted with R^{33};

4-($H_3CCH_2C(=O)$)-phenyl- substituted with R^{33};

4-($H_3CC(=O)$)-phenyl- substituted with R^{33};

4-($H_3CCH_2CH_2CH(OH)$)-phenyl- substituted with R^{33};

4-($(H_3C)_2CHCH(OH)$)-phenyl- substituted with R^{33};

4-($H_3CCH_2CH(OH)$)-phenyl- substituted with R^{33};

4-($H_3CCH(OH)$)-phenyl- substituted with R^{33};

4-(cyclopropyloxy)-phenyl- substituted with R^{33};

4-(cyclobutyloxy)-phenyl- substituted with R^{33}; or

4-(cyclopentyloxy)-phenyl- substituted with R^{33};

R^{12} is selected from

phenyl- substituted with 0-5 fluoro;

naphthyl- substituted with 0-3 R^{33};

2-($H_3CCH_2C(=O)$)-phenyl- substituted with R^{33};

2-($H_3CC(=O)$)-phenyl- substituted with R^{33};

2-($HC(=O)$)-phenyl- substituted with R^{33};

2-($H_3CCH(OH)$)-phenyl- substituted with R^{33};

2-($H_3CCH_2CH(OH)$)-phenyl- substituted with R^{33};

2-($HOCH_2$)-phenyl- substituted with R^{33};

2-($HOCH_2CH_2$)-phenyl- substituted with R^{33};

2-(H_3COCH_2)-phenyl- substituted with R^{33};

2-($H_3COCH_2CH_2$)-phenyl- substituted with R^{33};

2-($H_3CCH(OMe)$)-phenyl- substituted with R^{33};

2-($H_3COC(=O)$)-phenyl- substituted with R^{33};

2-($HOCH_2CH=CH$)-phenyl- substituted with R^{33};

2-((MeOC=O)CH=CH)-phenyl- substituted with R^{33};

2-(methyl)-phenyl- substituted with R^{33};

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2-(ethyl)-phenyl- substituted with R^{33};

2-(i-propyl)-phenyl- substituted with R^{33};

2-(F_3C)-phenyl- substituted with R^{33};

2-(NC)-phenyl- substituted with R^{33};

2-(H_3CO)-phenyl- substituted with R^{33};

2-(fluoro)-phenyl- substituted with R^{33};

2-(chloro)-phenyl- substituted with R^{33};

3-(NC)-phenyl- substituted with R^{33};

3-(H_3CO)-phenyl- substituted with R^{33};

3-(fluoro)-phenyl- substituted with R^{33};

3-(chloro)-phenyl- substituted with R^{33};

4-(NC)-phenyl- substituted with R^{33};

4-(fluoro)-phenyl- substituted with R^{33};

4-(chloro)-phenyl- substituted with R^{33};

4-(H_3CS)-phenyl- substituted with R^{33};

4-(H_3CO)-phenyl- substituted with R^{33};

4-(ethoxy)-phenyl- substituted with R^{33};

4-(i-propoxy)-phenyl- substituted with R^{33};

4-(i-butoxy)-phenyl- substituted with R^{33};

4-($H_3CCH_2CH_2C(=O)$)-phenyl- substituted with R^{33};

4-($(H_3C)_2CHC(=O)$)-phenyl- substituted with R^{33};

4-($H_3CCH_2C(=O)$)-phenyl- substituted with R^{33};

4-($H_3CC(=O)$)-phenyl- substituted with R^{33};

4-($H_3CCH_2CH_2CH(OH)$)-phenyl- substituted with R^{33};

4-($(H_3C)_2CHCH(OH)$)-phenyl- substituted with R^{33};

4-($H_3CCH_2CH(OH)$)-phenyl- substituted with R^{33};

4-($H_3CCH(OH)$)-phenyl- substituted with R^{33};

4-(cyclopropyloxy)-phenyl- substituted with R^{33};

4-(cyclobutyloxy)-phenyl- substituted with R^{33}; and

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4-(cyclopentyloxy)-phenyl- substituted with R^{33};

R^{13} is H, methyl, or ethyl;

alternatively, R^{12} and R^{13} join to form a 5- or 6-membered ring selected from pyrrolyl, pyrrolidinyl, imidazolyl, piperidinyl, piperizinyl, methylpiperizinyl,and morpholinyl;

alternatively, R^{12} and R^{13} when attached to N may be combined to form a 9- or 10-membered bicyclic heterocyclic ring system containing from 1-3 heteroatoms selected from the group consisting of N, O, and S; wherein said bicyclic heterocyclic ring system is selected from indolyl, indolinyl, indazolyl, benzimidazolyl, benzimidazolinyl, and benztriazolyl; wherein said bicyclic heterocyclic ring system is substituted with 0-1 R^{16};

R^{16}, at each occurrence, is independently selected from
H, OH, F, Cl, CN, NO_2, methyl, ethyl, methoxy, ethoxy, trifluoromethyl, and trifluoromethoxy; and

R^{31}, at each occurrence, is independently selected from
H, -OH, F, Cl, $-CF_3$, $-OCF_3$, methyl, ethyl,

methyl-C(=O)-, ethyl-C(=O)-, methoxy, ethoxy, methylthio-, ethylthio-, methyl-S(=O)-, ethyl-S(=O)-, methyl-SO_2-, and ethyl-SO_2-;

R^{33}, at each occurrence, is independently selected from

H, F, Cl, -CH_3, -OCH_3, -CF_3, -OCF_3, -CN, and -NO_2.

3. (Original) A compound of Claim 2 of Formula (I) wherein:

R^A is selected from

phenyl-NH-, (1-naphthyl)-NH-,
(2-naphthyl)-NH-, (2-[1,1'-biphenyl])-NH-,
(3-[1,1'-biphenyl])-NH-, (4-[1,1'-biphenyl])-NH-,
(2-F-phenyl)-NH-, (2-Cl-phenyl)-NH-,
(2-CF_3-phenyl)-NH-, (2-CH_3-phenyl)-NH-,
(2-OMe-phenyl)-NH-, (2-CN-phenyl)-NH-,
(2-OCF_3-phenyl)-NH-, (2-SMe-phenyl)-NH-,
(3-F-phenyl)-NH-, (3-Cl-phenyl)-NH-,
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(4-OCF_3-phenyl)-NH-, (4-SMe-phenyl)-NH-,
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(2,5-diCl-phenyl)-NH-, (2,6-diCl-phenyl)-NH-,
(3,4-diCl-phenyl)-NH-, (3,5-diCl-phenyl)-NH-,
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(2,3-diCF$_3$-phenyl)-NH-, (2,4-diCF$_3$-phenyl)-NH-,
(2,5-diCF$_3$-phenyl)-NH-, (2,6-diCF$_3$-phenyl)-NH-,
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A4 (2-F-3-OMe-phenyl)-NH-, (2-F-4-OMe-phenyl)-NH-,
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(2-OMe-5-CH_3-phenyl)-NH-, (2-OMe-6-CH_3-phenyl)-NH-,
A4 (2-OMe-3-CF_3-phenyl)-NH-, (2-OMe-4-CF_3-phenyl)-NH-,
(2-OMe-5-CF_3-phenyl)-NH-, (2-OMe-6-CF_3-phenyl)-NH-

(3-CF_3-4-Cl-phenyl)-NH-, (3-CF_3-4-C(O)CH_3-phenyl)-NH-, (2,3,5-triCl-phenyl)-NH-, (3-CH_3-4-CO_2Me-phenyl)-NH-, and
(3-CHO-4-OMe-phenyl)-NH-;

R^B is selected from
2-chlorophenyl, 2-fluorophenyl, 2-bromophenyl,
2-cyanophenyl, 2-methylphenyl, 2-trifluoromethylphenyl,
2-methoxyphenyl, 2-trifluoromethoxyphenyl,

3-chlorophenyl, 3-fluorophenyl, 3-bromophenyl,
3-cyanophenyl, 3-methylphenyl, 3-ethylphenyl,
3-propylphenyl, 3-isopropylphenyl, 3-butylphenyl,
3-trifluoromethylphenyl, 3-methoxyphenyl,
3-isopropoxyphenyl, 3-trifluoromethoxyphenyl,
3-thiomethoxyphenyl,

4-chlorophenyl, 4-fluorophenyl, 4-bromophenyl,
4-cyanophenyl, 4-methylphenyl, 4-ethylphenyl,
4-propylphenyl, 4-isopropylphenyl, 4-butylphenyl,
4-trifluoromethylphenyl, 4-methoxyphenyl,
4-isopropoxyphenyl, 4-trifluoromethoxyphenyl,
4-thiomethoxyphenyl,

2,3-dichlorophenyl, 2,3-difluorophenyl,
2,3-dimethylphenyl, 2,3-ditrifluoromethylphenyl,
2,3-dimethoxyphenyl, 2,3-ditrifluoromethoxyphenyl,

2,4-dichlorophenyl, 2,4-difluorophenyl,
2,4-dimethylphenyl, 2,4-ditrifluoromethylphenyl,
2,4-dimethoxyphenyl, 2,4-ditrifluoromethoxyphenyl,

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2,5-dichlorophenyl, 2,5-difluorophenyl,
2,5-dimethylphenyl, 2,5-ditrifluoromethylphenyl,
2,5-dimethoxyphenyl, 2,5-ditrifluoromethoxyphenyl,

2,6-dichlorophenyl, 2,6-difluorophenyl,
2,6-dimethylphenyl, 2,6-ditrifluoromethylphenyl,
2,6-dimethoxyphenyl, 2,6-ditrifluoromethoxyphenyl,

3,4-dichlorophenyl, 3,4-difluorophenyl,
3,4-dimethylphenyl, 3,4-ditrifluoromethylphenyl,
3,4-dimethoxyphenyl, 3,4-ditrifluoromethoxyphenyl,

2,4,6-trichlorophenyl, 2,4,6-trifluorophenyl,
2,4,6-trimethylphenyl, 2,4,6-tritrifluoromethylphenyl,
2,4,6-trimethoxyphenyl, 2,4,6-tritrifluoromethoxyphenyl,

2-chloro-4-CF_3-phenyl, 2-fluoro-3-chloro-phenyl,
2-chloro-4-CF_3-phenyl, 2-chloro-4-methoxy-phenyl,
2-methoxy-4-isopropyl-phenyl, 2-CF_3-4-methoxy-phenyl,

2-methyl-4-methoxy-5-fluoro-phenyl,
2-methyl-4-methoxy-phenyl, 2-chloro-4-CF_3O-phenyl,
2,4,5-trimethyl-phenyl, 2-methyl-4-chloro-phenyl,

4-acetylphenyl, 3-acetamidophenyl, 2-naphthyl;

2-Me-5-F-phenyl, 2-F-5-Me-phenyl, 2-MeO-5-F-phenyl,
2-Me-3-Cl-phenyl, 3-NO_2-phenyl, 2-NO_2-phenyl,
2-Cl-3-Me-phenyl, 2-Me-4-EtO-phenyl, 2-Me-4-F-phenyl,
2-Cl-6-F-phenyl, 2-Cl-4-(CHF_2)O-phenyl,
2,4-diMeO-6-F-phenyl, 2-CF_3-6-F-phenyl,
2-MeS-phenyl, 2,6-diCl-4-MeO-phenyl,
2,3,4-triF-phenyl, 2,6-diF-4-Cl-phenyl,
2,3,4,6-tetraF-phenyl, 2,3,4,5,6-pentaF-phenyl,
2-CF_3-4-EtO-phenyl, 2-CF_3-4-iPrO-phenyl,
2-CF_3-4-Cl-phenyl, 2-CF_3-4-F-phenyl, 2-Cl-4-EtO-phenyl,
2-Cl-4-iPrO-phenyl, 2-Et-4-MeO-phenyl,
2-CHO-4-MeO-phenyl, 2-CH_3CH(OH)-4-MeO-phenyl,
2-CH_3CH(OH)-4-F-phenyl, 2-CH_3CH(OH)-4-Cl-phenyl,
2-CH_3CH(OH)-4-Me-phenyl, 2-CH_3CH(OMe)-4-MeO-phenyl,
2-CH_3C(=O)-4-MeO-phenyl, 2-CH_3C(=O)-4-F-phenyl,
2-CH_3C(=O)-4-Cl-phenyl, 2-CH_3C(=O)-4-Me-phenyl,
2-H_2C(OH)-4-MeO-phenyl, 2-H_2C(OMe)-4-MeO-phenyl,
2-H_3CCH_2CH(OH)-4-MeO-phenyl, 2-H_3CCH_2C(=O)-4-MeO-phenyl,

2-$CH_3CO_2CH_2CH_2$-4-MeO-phenyl,
(Z)-2-$HOCH_2CH$=CH-4-MeO-phenyl,
(E)-2-$HOCH_2CH$=CH-4-MeO-phenyl,
(Z)-2-CH_3CO_2CH=CH-4-MeO-phenyl,
(E)-2-CH_3CO_2CH=CH-4-MeO-phenyl,
2-$CH_3OCH_2CH_2$-4-MeO-phenyl,
2-F-4-MeO-phenyl, 2-Cl-4-F-phenyl,
cyclohexyl, cyclopentyl, cyclohexylmethyl,

benzyl, 2-F-benzyl, 3-F-benzyl, 4-F-benzyl,
3-MeO-benzyl, 3-OH-benzyl, 2-MeO-benzyl,
2-OH-benzyl, 2-MeOC(=O)-3-MeO-phenyl,
2-Me-4-CN-phenyl, 2-Me-3-CN-phenyl,
2-Me-4-MeS-phenyl, 2-CF_3-4-CN-phenyl,
2-CHO-phenyl, 3-CHO-phenyl, 2-$HOCH_2$-phenyl,
3-$HOCH_2$-phenyl, 3-$MeOCH_2$-phenyl,
3-Me_2NCH_2-phenyl, 3-CN-4-F-phenyl,
2-Me-4-H_2NCO-phenyl, 2-Me-4-MeOC(=O)-phenyl,
3-H_2NCO-4-F-phenyl, 2-Me_2NCH_2-4-MeO-phenyl-,
2-Me-4-CH_3C(=O)-phenyl,

phenyl-NH-, (1-naphthyl)-NH-,
(2-naphthyl)-NH-, (2-[1,1'-biphenyl])-NH-,
(3-[1,1'-biphenyl])-NH-, (4-[1,1'-biphenyl])-NH-,
(2-F-phenyl)-NH-, (2-Cl-phenyl)-NH-,
(2-CF_3-phenyl)-NH-, (2-CH_3-phenyl)-NH-,
(2-OMe-phenyl)-NH-, (2-CN-phenyl)-NH-,
(2-OCF_3-phenyl)-NH-, (2-SMe-phenyl)-NH-,
(3-F-phenyl)-NH-, (3-Cl-phenyl)-NH-,
(3- CF_3-phenyl)-NH-, (3-CH_3-phenyl)-NH-,
(3-OMe-phenyl)-NH-, (3-CN-phenyl)-NH-,
(3-OCF_3-phenyl)-NH-, (3-SMe-phenyl)-NH-,
(4-F-phenyl)-NH-, (4-Cl-phenyl)-NH-,
(4-CF_3-phenyl)-NH-, (4-CH_3-phenyl)-NH-,
(4-OMe-phenyl)-NH-, (4-CN-phenyl)-NH-,
(4-OCF_3-phenyl)-NH-, (4-SMe-phenyl)-NH-,
(2,3-diCl-phenyl)-NH-, (2,4-diCl-phenyl)-NH-,
(2,5-diCl-phenyl)-NH-, (2,6-diCl-phenyl)-NH-,
(3,4-diCl-phenyl)-NH-, (3,5-diCl-phenyl)-NH-,
(2,3-diF-phenyl)-NH-, (2,4-diF-phenyl)-NH-,
(2,5-diF-phenyl)-NH-, (2,6-diF-phenyl)-NH-,
(3,4-diF-phenyl)-NH-, (3,5-diF-phenyl)-NH-,

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(2,3-diCH_3-phenyl)-NH-, (2,4-diCH_3-phenyl)-NH-,
(2,5-diCH_3-phenyl)-NH-, (2,6-diCH_3-phenyl)-NH-,
(3,4-diCH_3-phenyl)-NH-, (3,5-diCH_3-phenyl)-NH-,
(2,3-diCF_3-phenyl)-NH-, (2,4-diCF_3-phenyl)-NH-,
(2,5-diCF_3-phenyl)-NH-, (2,6-diCF_3-phenyl)-NH-,
(3,4-diCF_3-phenyl)-NH-, (3,5-diCF_3-phenyl)-NH-,
(2,3-diOMe-phenyl)-NH-, (2,4-diOMe-phenyl)-NH-,
(2,5-diOMe-phenyl)-NH-, (2,6-diOMe-phenyl)-NH-,
(3,4-diOMe-phenyl)-NH-, (3,5-diOMe-phenyl)-NH-,

(2-F-3-Cl-phenyl)-NH-, (2-F-4-Cl-phenyl)-NH-,
(2-F-5-Cl-phenyl)-NH-, (2-F-6-Cl-phenyl)-NH-,
(2-F-3-CH_3-phenyl)-NH-, (2-F-4-CH_3-phenyl)-NH-,
(2-F-5-CH_3-phenyl)-NH-, (2-F-6-CH_3-phenyl)-NH-,
(2-F-3-CF_3-phenyl)-NH-, (2-F-4-CF_3-phenyl)-NH-,
A4 (2-F-5-CF_3-phenyl)-NH-, (2-F-6-CF_3-phenyl)-NH-,
(2-F-3-OMe-phenyl)-NH-, (2-F-4-OMe-phenyl)-NH-,
(2-F-5-OMe-phenyl)-NH-, (2-F-6-OMe-phenyl)-NH-,

(2-Cl-3-F-phenyl)-NH-, (2-Cl-4-F-phenyl)-NH-,
(2-Cl-5-F-phenyl)-NH-, (2-Cl-6-F-phenyl)-NH-,
(2-Cl-3-CH_3-phenyl)-NH-, (2-Cl-4-CH_3-phenyl)-NH-,
(2-Cl-5-CH_3-phenyl)-NH-, (2-Cl-6-CH_3-phenyl)-NH-,
(2-Cl-3-CF_3-phenyl)-NH-, (2-Cl-4-CF_3-phenyl)-NH-,
(2-Cl-5-CF_3-phenyl)-NH-, (2-Cl-6-CF_3-phenyl)-NH-,
(2-Cl-3-OMe-phenyl)-NH-, (2-Cl-4-OMe-phenyl)-NH-,
(2-Cl-5-OMe-phenyl)-NH-, (2-Cl-6-OMe-phenyl)-NH-,

(2-CH_3-3-F-phenyl)-NH-, (2-CH_3-4-F-phenyl)-NH-,
(2-CH_3-5-F-phenyl)-NH-, (2-CH_3-6-F-phenyl)-NH-,
(2-CH_3-3-Cl-phenyl)-NH-, (2-CH_3-4-Cl-phenyl)-NH-,
(2-CH_3-5-Cl-phenyl)-NH-, (2-CH_3-6-Cl-phenyl)-NH-,
(2-CH_3-3-CF_3-phenyl)-NH-, (2-CH_3-4-CF_3-phenyl)-NH-,

(2-CH_3-5-CF_3-phenyl)-NH-, (2-CH_3-6-CF_3-phenyl)-NH-,
(2-CH_3-3-OMe-phenyl)-NH-, (2-CH_3-4-OMe-phenyl)-NH-,
(2-CH_3-5-OMe-phenyl)-NH-, (2-CH_3-6-OMe-phenyl)-NH-,

(2-CF_3-3-F-phenyl)-NH-, (2-CF_3-4-F-phenyl)-NH-,
(2-CF_3-5-F-phenyl)-NH-, (2-CF_3-6-F-phenyl)-NH-,
(2-CF_3-3-Cl-phenyl)-NH-, (2-CF_3-4-Cl-phenyl)-NH-,
(2-CF_3-5-Cl-phenyl)-NH-, (2-CF_3-6-Cl-phenyl)-NH-,
(2-CF_3-3-CH_3-phenyl)-NH-, (2-CF_3-4-CH_3-phenyl)-NH-,
(2-CH_3-5-CF_3-phenyl)-NH-, (2-CF_3-6-CH_3-phenyl)-NH-,
(2-CF_3-3-OMe-phenyl)-NH-, (2-CF_3-4-OMe-phenyl)-NH-,
(2-CF_3-5-OMe-phenyl)-NH-, (2-CF_3-6-OMe-phenyl)-NH-,

(2-OMe-3-F-phenyl)-NH-, (2-OMe-4-F-phenyl)-NH-,
(2-OMe-5-F-phenyl)-NH-, (2-OMe-6-F-phenyl)-NH-,
(2-OMe-3-Cl-phenyl)-NH-, (2-OMe-4-Cl-phenyl)-NH-,
(2-OMe-5-Cl-phenyl)-NH-, (2-OMe-6-Cl-phenyl)-NH-,
(2-OMe-3-CH_3-phenyl)-NH-, (2-OMe-4-CH_3-phenyl)-NH-,
(2-OMe-5-CH_3-phenyl)-NH-, (2-OMe-6-CH_3-phenyl)-NH-,
(2-OMe-3-CF_3-phenyl)-NH-, (2-OMe-4-CF_3-phenyl)-NH-,
(2-OMe-5-CF_3-phenyl)-NH-, (2-OMe-6-CF_3-phenyl)-NH-

(3-CF_3-4-Cl-phenyl)-NH-, (3-CF_3-4-C(O)CH_3-phenyl)-NH-, (2,3,5-triCl-phenyl)-NH-, (3-CH_3-4-CO_2Me-phenyl)-NH-, and
(3-CHO-4-OMe-phenyl)-NH-; and

R^{31}, at each occurrence, is independently selected from
H, -OH, F, Cl, -CF_3, -OCF_3, methyl, methyl-C(=O)-, methoxy, methylthio-, methyl-S(=O)-, and methyl-SO_2-.

4. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.

5. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 2, or a pharmaceutically acceptable salt thereof.

6. (Original) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 3, or a pharmaceutically acceptable salt thereof.

7. (Original) A method for treating a human suffering from a disorder associated with 5HT2C receptor modulation comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.

8. (Original) A method for treating a human suffering from a disorder associated with 5HT2C receptor modulation comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 2, or a pharmaceutically acceptable salt thereof.

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9. (Original) A method for treating a human suffering from a disorder associated with 5HT2C receptor modulation comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 3, or a pharmaceutically acceptable salt thereof.

10. (Original) A method of Claim 7 for treating a human suffering from a disorder associated with 5HT2C receptor modulation wherein the compound is a 5HT2C agonist.

11. (Original) A method of Claim 8 for treating a human suffering from a disorder associated with 5HT2C receptor modulation wherein the compound is a 5HT2C agonist.

12. (Original) A method of Claim 9 for treating a human suffering from a disorder associated with 5HT2C receptor modulation wherein the compound is a 5HT2C agonist.

13. (Original) A method for treating obesity comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.

14. (Original) A method for treating obesity comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 2, or a pharmaceutically acceptable salt thereof.

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15. (Original) A method for treating obesity comprising administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 3, or a pharmaceutically acceptable salt thereof.